CLAIM LISTING

1. (Previously presented) A compound of formula (Ia)

$$R^{2}$$
 R^{3}
 R^{4}
 $(CH_{2})_{n}$
 R^{5}
 $(O)_{m}$
 Ar
 (Ia)

wherein R^1 , R^2 , R^3 , and R^4 independently of each other represent hydrogen, halogen, perhalomethyl, hydroxy, nitro, cyano, formyl, or C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, C_{1-12} -alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxy C_{1-12} -alkyl, amino, acylamino, C_{1-12} -alkylamino, arylamino, aralkylamino, amino C_{1-12} -alkyl, C_{1-12} -alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, C_{1-12} -alkoxy C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, aralkoxy C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, - COR^{11} , or - SO_2R^{12} , wherein R^{11} and R^{12} independently of each other are selected from hydroxy, halogen, perhalomethyl, C_{1-6} -alkoxy or amino optionally substituted with one or more C_{1-6} -alkyl, perhalomethyl or aryl, optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^1 and R^2 , R^2 and R^3 and/or R^3 and R^4 may form a cyclic ring containing from 5 to 7 carbon atoms optionally substituted with one or more C_{1-6} -alkyl;

ring A fused to the ring containing X and N represents a 5-6 membered cyclic ring, optionally substituted with one or more hydrogen, halogen, perhalomethyl, hydroxy or C₁₋₇-alkyl, C₂₋₇-alkynyl, C₁₋₇-alkoxy or aryl,

X is a -(CHR 9)-CH $_2$ -, -CH=CH-, -(NR 9)-CH $_2$ -, -(CHR 9)-CH=CH-, -(CHR 9)-CH $_2$ -CH $_2$ -, -CH=(CR 9)-, -(CO)-(CHR 9)-, wherein R 9 is hydrogen, halogen, hydroxy, nitro, cyano, formyl, C $_{1\cdot12}$ -alkyl, C $_{1\cdot12}$ -alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxyalkyl, amino, acylamino, C $_{1\cdot12}$ -alkylamino, arylamino, aralkylamino, aminoC $_{1\cdot12}$ -alkyl, C $_{1\cdot12}$ -alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, C $_{1\cdot12}$ -alkoxyC $_{1\cdot12}$ -alkyl, aralkoxyC $_{1\cdot12}$ -alkyl, C $_{1\cdot12}$ -alkyl, thio, thioC $_{1\cdot12}$ -alkyl, C $_{1\cdot12}$ -alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, -COR $_{13}$, or -SO $_2$ R $_1^{14}$, wherein R $_1^{13}$ and R $_1^{14}$ independently of each other are selected from hydroxy, halogen, C $_{1\cdot6}$ -alkoxy, amino optionally substituted with one or more C $_{1\cdot6}$ -alkyl, perhalomethyl or aryl;

Ar represents arylene or heteroarylene, optionally substituted with one or more C_{1-6} -alkyl or aryl;

 R^5 represents hydrogen, hydroxy, halogen, C_{1-12} -alkoxy, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl or aralkyl, optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^5 forms a bond together with R^6 , R^6 represents hydrogen, hydroxy, halogen, C_{1-12} -alkoxy, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, acyl or aralkyl; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^6 forms a bond together with R^5 , R^7 represents hydrogen, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, aryl, aralkyl, C_{1-12} -alkoxy C_{1-12} -alkoxycarbonyl, aryloxycarbonyl, C_{1-12} -alkylaminocarbonyl, arylaminocarbonyl, acyl, heterocyclyl, heteroaryl or heteroaralkyl groups; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; R^8 represents hydrogen, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, aryl, aralkyl, heterocyclyl, heteroaryl or heteroaralkyl groups; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano;

Y represents oxygen, sulphur or NR^{10} , where R^{10} represents hydrogen, C_{1-12} -alkyl, aryl, hydroxy C_{1-12} -alkyl or aralkyl groups or when Y is NR^{10} , R^8 and R^{10} may form a 5 or 6 membered nitrogen containing ring, optionally substituted with one or more C_{1-6} -alkyl; n is an integer ranging from 1 to 4 and m is an integer ranging from 0 to 1;

or a pharmaceutically acceptable salt thereof.

2. (Original) A compound according to claim 1 wherein R¹, R², R³, and R⁴ independently of each other represent hydrogen, halogen, perhalomethyl, hydroxy, cyano, or C₁₋₇-alkyl, C₄₋₇-alkenynyl, C₂₋₇-alkenyl, C₂₋₇-alkynyl, C₁₋₇-alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxyC₁₋₇-alkyl, amino, acylamino, C₁₋₇-alkylamino, arylamino, aralkylamino, aminoC₁₋₇-alkyl, C₁₋₇-alkoxyC₁₋₇-alkyl, aryloxyC₁₋₇-alkyl, aralkoxyC₁₋₇-alkyl, C₁₋₇-alkyl, C₁₋₇-alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, -COR¹¹, or -SO₂R¹², wherein R¹¹ and R¹² independently of each other are selected from hydroxy, perhalomethyl or amino optionally substituted with one or more C₁₋₆-alkyl, perhalomethyl or aryl; optionally substituted with one or more halogen, perhalomethyl, hydroxy or cyano; or R¹ and R², R² and R³ and/or R³ and R⁴ may form a cyclic ring containing from 5 to 7 carbon atoms optionally substituted with one or more C₁₋₆-alkyl.

- 3. (Cancelled)
- 4. (Cancelled)
- 5. (Cancelled)
- 6. (Cancelled)

7. (Previously presented) A compound according to claim 1 wherein ring A fused to the ring containing X and N represents a 5-6 membered cyclic ring, optionally substituted with one or more hydrogen, halogen, perhalomethyl, hydroxy, cyano, or C₁₋₇-alkyl, C₄₋₇-alkenynyl, C₂₋₇-alkenyl, C₂₋₇-alkynyl, C₁₋₇-alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxyC₁₋₇-alkyl, amino, acylamino, C₁₋₇-alkylamino, arylamino, aralkylamino, aminoC₁₋₇-alkyl, C₁₋₇-alkoxyC₁₋₇-alkyl, aralkoxyC₁₋₇-alkyl, C₁₋₇-alkylthio, thioC₁₋₇-alkyl, C₁₋₇-alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, -COR¹¹, or -SO₂R¹², wherein R¹¹ and R¹² independently of each other are selected from hydroxy, perhalomethyl or amino optionally

substituted with one or more C_{1-6} -alkyl, perhalomethyl or aryl; optionally substituted with one or more halogen, perhalomethyl, hydroxy or cyano.

- 8. (Cancelled)
- 9. (Cancelled)
- 10. (Cancelled)
- 11. (Cancelled)
- 12. (Cancelled)
- 13. (Cancelled)
- 14. (Cancelled)
- 15. (Cancelled)
- 16. (Cancelled)
- 17. (Previously presented) A compound according to claim 1 wherein Ar represents arylene or heteroarylene;

R⁵ represents hydrogen, hydroxy, halogen; or R⁵ forms a bond together with R⁶,

R⁶ represents hydrogen, hydroxy, halogen; or R⁶ forms a bond together with R⁵,

 R^7 represents hydrogen, C_{1-7} -alkyl, C_{2-7} -alkenyl, C_{2-7} -alkynyl, aryl, aralkyl, C_{1-7} -alkyl, C_{1-7} -alkylaminocarbonyl, arylaminocarbonyl, acyl, heterocyclyl, heteroaryl or

heteroaralkyl groups;

 R^8 represents hydrogen, C_{1-7} -alkyl, C_{2-7} -alkenyl, C_{2-7} -alkynyl;

Y represents oxygen or sulphur;

n is an integer ranging from 2 to 3 and m is 1.

- 18. (Cancelled)
- 19. (Cancelled)
- 20. (Cancelled)
- 21. (Cancelled)
- 22. (Cancelled)
- 23. (Cancelled)
- 24. (Cancelled)

- 25. (Cancelled)
- 26. (Cancelled)
- 27. (Cancelled)
- 28. (Cancelled)
- 29. (Cancelled)
- 30. (Cancelled)
- 31. (Cancelled)
- 32. (Cancelled)
- 33. (Cancelled)
- 34. (Cancelled)
- 35. (Cancelled)
- 36. (Cancelled)
- 37. (Cancelled)
- 38. (Cancelled)
- 39. (Cancelled)
- 40. (Cancelled)
- 41. (Cancelled)
- 42. (Cancelled)
- 43. (Cancelled)
- 44. (Cancelled)
- 45. (Previously presented) The compound according to claim 1 which is
- 3-{4-[2-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-2-ethoxy-propionic acid,
- Ethyl-3- $\{4-[2-(10,11-dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl\}-2-ethoxy-propionionate,$
- 3-{4-[2-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-2-methoxy-propionic acid,
- 3-{4-[2-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-2-propoxy-propionic acid,
- 3-{4-[2-(10,11-Dihydro-dibenzo[*b,f*]azepin-5-yl)-ethoxy]-phenyl}-2-benzyloxy-propionic acid,
- 3-{4-[3-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-propoxy]-phenyl}-2-ethoxy-propionic acid,

- 3-{4-[3-(10,11-Dihydro-dibenzo[*b,f*]azepin-5-yl)-propoxy]-phenyl}-2-methoxy-propionic acid,
- 3-{4-[3-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl}-2-ethoxy-propionic acid,
- 3-{4-[3-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl}-2-methoxy-propionic acid,
- 3-{4-[2-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-methoxy]-phenyl}-2-ethoxy-propionic acid,
- 2-Ethoxy-3-{4-[2-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid,
- 2-Methoxy-3-{4-[2-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid,
- 2-Propoxy-3-{4-[2-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-ethoxy]-phenyl}-propionic acid,
- 2-Benzyloxy-3-{4-[2-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-ethoxy]-phenyl}-propionic acid,
- 2-Ethoxy-3-{4-[1-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-methoxy]-phenyl}-propionic acid,
- 2-Ethoxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propoxy]-phenyl}-propionic acid,
- 2-Methoxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propoxy]-phenyl}-propionic acid,
- 2-Propoxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propoxy]-phenyl}-propionic acid,
- 2-Benzyloxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propoxy]-phenyl}-propionic acid,
- 2-Ethoxy-3- $\{4-[3-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl\}-propionic acid.$
- 2-Methoxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propyl]-phenyl}-propionic acid,
- 2-Propoxy-3-{4-[3-(10-oxo-10,11-dihydro-dibenzo[*b,f*]azepin-5-yl)-propyl]-phenyl}-propionic acid,

- 2-Benzyloxy-3- $\{4-[3-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl\}$ propionic acid, 2-Ethoxy-3-{4-[2-(10-methoxy-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid, 2-Methoxy-3-{4-[2-(10-methoxy-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid, 2-Propoxy-3-{4-[2-(10-methoxy-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid. 2-Ethoxy-3-{4-[1-(10-methoxy-dibenzo[b,f]azepin-5-yl)-methoxy]-phenyl}-propionic acid. 2-Benzyloxy-3-{4-[2-(10-methoxy-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid, 2-Ethoxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propoxy]-phenyl}-propionic acid, 2-Methoxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propoxy]-phenyl}-propionic acid, 2-Benzyloxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propoxy]-phenyl}-propionic acid. 2-Ethoxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl}-propionic acid, 2-Methoxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl}-propionic acid, 2-Benzyloxy-3-{4-[3-(10-methoxy-dibenzo[b,f]azepin-5-yl)-propyl]-phenyl}-propionic acid, 3-(4-(2-(Dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-ethoxy-propionic acid. Ethyl-3-(4-(2-(dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-ethoxy-propionionate,3-(4-(2-(Dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-methoxy-propionic acid, 3-(4-(2-(Dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-propoxy-propionic acid, 3-(4-(2-(Dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-benzyloxy-propionic acid. 3-(4-(1-(Dibenzo[b,f]azepin-5-yl)-methoxy)-phenyl)-2-ethoxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propoxy)-phenyl)-2-ethoxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propoxy)-phenyl)-2-propoxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propoxy)-phenyl)-2-benzyloxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propyl)-phenyl)-2-ethoxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propyl)-phenyl)-2-propoxy-propionic acid, 3-(4-(3-(Dibenzo[b,f]azepin-5-yl)-propyl)-phenyl)-2-benzyloxy-propionic acid, or a pharmaceutically acceptable salt thereof.
- 46. (Previously presented) The compound according to claim 1 which is 3-{4-[2-(10,11-Dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-2-ethoxy-propionic acid,

Ethyl-3-(4-(2-(10,11-dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy)-phenyl)-2-ethoxy-propionionate,

- 2-Ethoxy-3-{4-[2-(10-oxo-10,11-dihydro-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid.
- 2-Ethoxy-3-{4-[2-(10-methoxy-dibenzo[b,f]azepin-5-yl)-ethoxy]-phenyl}-propionic acid, or a pharmaceutically acceptable salt thereof.
- 47. (Previously presented) A pharmaceutical composition comprising, as an active ingredient, a compound according to claim 1 or a pharmaceutically acceptable salt thereof together with a pharmaceutically acceptable carrier or diluent.
- 48. (Cancelled)
- 49. (Cancelled)
- 52. (Cancelled)
- 51. (Cancelled)
- 52. (Cancelled)
- 53. (Cancelled)
- 54. (Currently amended) A method for the treatment of conditions mediated by nuclear receptors, in particular the Peroxisome Proliferator-Activated Receptors (PPAR), the method comprising administering to a subject in need thereof an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof.
- 55. (Previously presented) A method for the treatment of diabetes or obesity, the method comprising administering to a subject in need thereof an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof.
- 56. (Cancelled)
- 57. (Cancelled)
- 58. (Cancelled)
- 59. (Cancelled)
- 60. (Cancelled)